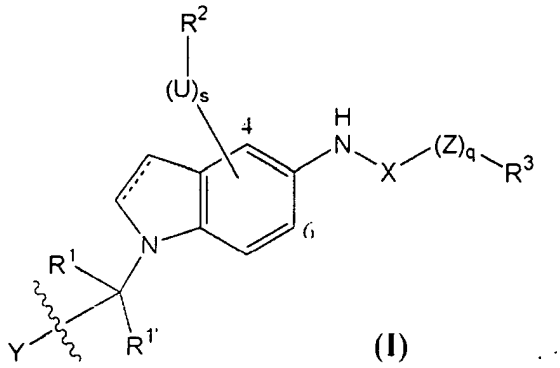
**Amendments to the Claims:**

This listing of claims will replace all prior versions, and listings, of claims in the application.

**Listing of Claims:**

1. (Currently amended) A compound of formula I:




**(I)**

wherein:

the dotted line represents an optional bond;

$\mathbf{R^1}$ and $\mathbf{R^{1'}}$ are each independently selected from the group consisting of hydrogen, $C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yl, $C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, hydroxy-$C_{1-6}$-alk(en/yn)yl, hydroxy-$C_{3-8}$-cycloalk(en)yl, hydroxy-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, halo-$C_{1-6}$-alk(en/yn)yl, halo-$C_{3-8}$-cycloalk(en)yl, halo-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, cyano-$C_{1-6}$-alk(en/yn)yl, cyano-$C_{3-8}$-cycloalk(en)yl and cyano-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl; or $\mathbf{R^1}$ and $\mathbf{R^{1'}}$ taken together with the carbon atom to which they are attached form a 3-8 membered saturated or unsaturated ring [[which]]that optionally contains 1 or 2 heteroatoms;

s is 0 or 1;

**U** is O, $N\mathbf{R^{11}}$, S, $SO_2$, $SO_2N\mathbf{R^{11}}$, CO-O or CO-$N\mathbf{R^{11}}$; wherein $\mathbf{R^{11}}$ is selected from the group consisting of hydrogen, $C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yl, and $C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl; or $\mathbf{R^2}$ and $\mathbf{R^{11}}$ taken together with the nitrogen atom to which they are attached

form a 4-8 membered saturated or unsaturated ring [[which]]that optionally contains 1, 2 or 3 further heteroatoms;

$\mathbf{R^2}$ is selected from the group consisting of hydrogen, $C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yl, $C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, Ar, Ar-$C_{1-6}$-alk(en/yn)yl, Ar-$C_{3-8}$-cycloalk(en)yl, Ar-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, acyl, hydroxy-$C_{1-6}$-alk(en/yn)yl, hydroxy-$C_{3-8}$-cycloalk(en)yl, hydroxy-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, halogen, halo-$C_{1-6}$-alk(en/yn)yl, halo-$C_{3-8}$-cycloalk(en)yl, halo-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, cyano, cyano-$C_{1-6}$-alk(en/yn)yl, cyano-$C_{3-8}$-cycloalk(en)yl, cyano-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, -$NO_2$, $N\mathbf{R^{10}R^{10'}}$-$C_{1-6}$-alk(en/yn)yl, $N\mathbf{R^{10}R^{10'}}$-$C_{3-8}$-cycloalk(en)yl and $N\mathbf{R^{10}R^{10'}}$-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl; wherein

> $\mathbf{R^{10}}$ and $\mathbf{R^{10'}}$ are each independently selected from the group consisting of hydrogen, $C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yl, $C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, hydroxy-$C_{1-6}$-alk(en/yn)yl, hydroxy-$C_{3-8}$-cycloalk(en)yl, hydroxy-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, halo-$C_{1-6}$-alk(en/yn)yl, halo-$C_{3-8}$-cycloalk(en)yl, halo-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, cyano-$C_{1-6}$-alk(en/yn)yl, cyano-$C_{3-8}$-cycloalk(en)yl and cyano-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl[[,]]; or
>
> $\mathbf{R^{10}}$ and $\mathbf{R^{10'}}$ taken together with the nitrogen atom to which they are attached form a 4-8 membered saturated or unsaturated ring [[which]]that optionally contains 1, 2 or 3 further heteroatoms;

with the proviso that when $\mathbf{R^2}$ is $NO_2$, halogen or cyano, then **s** is 0; and

with the proviso that when $\mathbf{R^2}$ is a hydrogen atom or acyl and s is 1, then **U** is $N\mathbf{R^{11}}$, O or S;

wherein the group $-(\mathbf{U})_s-\mathbf{R^2}$ is linked to position 4 or 6 of the indole or indoline;

**q** is 0 or 1;

**Z** is O or S;

**X** is CO or $SO_2$; with the proviso that **q** is 0 when **X** is $SO_2$;

$\mathbf{R^3}$ is selected from the group consisting of $C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yl, heterocycloalk(en)yl, $C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, $C_{1-6}$-alk(en/yn)yl-$C_{3-8}$-cycloalk(en)yl, $C_{1-6}$-alk(en/yn)yl-heterocycloalk(en)yl, Ar, Ar-$C_{1-6}$-alk(en/yn)yl, Ar-$C_{3-8}$-cycloalk(en)yl, Ar-heterocycloalk(en)yl, Ar-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, Ar-$C_{1-6}$-alk(en/yn)yl-$C_{3-8}$-cycloalk(en)yl, Ar-$C_{1-6}$-alk(en/yn)yl-heterocycloalk(en)yl, $C_{1-6}$-alk(en/yn)yloxy-$C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yloxy-$C_{1-6}$-alk(en/yn)yl, $C_{1-6}$-alk(en/yn)yloxy-$C_{3-8}$-cycloalk(en)yl, $C_{1-6}$-alk(en/yn)yloxy-heterocycloalk(en)yl, Ar-oxy-$C_{1-6}$-alk(en/yn)yl, Ar-$C_{1-6}$-alk(en/yn)yloxy-$C_{1-6}$-alk(en/yn)yl, $C_{1-6}$-alk(en/yn)yloxy-carbonyl-$C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yloxy-carbonyl-$C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yloxy-carbonyl-$C_{1-6}$-alk(en/yn)yl, hydroxy-$C_{1-6}$-alk(en/yn)yl, hydroxy-$C_{3-8}$-cycloalk(en)yl, hydroxy-heterocycloalk(en)yl, hydroxy-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, hydroxy-$C_{1-6}$-alk(en/yn)yl-$C_{3-8}$-cycloalk(en)yl, hydroxy-$C_{1-6}$-alk(en/yn)yl-heterocycloalk(en)yl, halo-$C_{1-6}$-alk(en/yn)yl, halo-$C_{3-8}$-cycloalk(en)yl, halo-heterocycloalk(en)yl, halo-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, halo-$C_{1-6}$-alk(en/yn)yl-$C_{3-8}$-cycloalk(en)yl, halo-$C_{1-6}$-alk(en/yn)yl-heterocycloalk(en)yl, halo-$C_{1-6}$-alk(en/yn)yl-Ar, halo-$C_{3-8}$-cycloalk(en)yl-Ar, halo-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl-Ar, halo-$C_{1-6}$-alk(en/yn)yl-$C_{3-8}$-cycloalk(en)yl-Ar, cyano-$C_{1-6}$-alk(en/yn)yl, cyano-$C_{3-8}$-cycloalk(en)yl, cyano-heterocycloalk(en)yl, cyano-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, cyano-$C_{1-6}$-alk(en/yn)yl-$C_{3-8}$-cycloalk(en)yl, cyano-$C_{1-6}$-alk(en/yn)yl-heterocycloalk(en)yl, acyl-$C_{1-6}$-alk(en/yn)yl, acyl-$C_{3-8}$-cycloalk(en)yl, acyl-heterocycloalk(en)yl, acyl-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, acyl-$C_{1-6}$-alk(en/yn)yl-$C_{3-8}$-cycloalk(en)yl, acyl-$C_{1-6}$-alk(en/yn)yl-heterocycloalk(en)yl, [[ and ]]-$N\mathbf{R^{12}R^{12'}}$, optionally substituted $N\mathbf{R^{12}R^{12'}}$-$C_{1-6}$-alk(en/yn)yl, optionally substituted $N\mathbf{R^{12}R^{12'}}$-$C_{3-8}$-cycloalk(en)yl, and optionally substituted $N\mathbf{R^{12}R^{12'}}$-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl; wherein

$\mathbf{R^{12}}$ and $\mathbf{R^{12'}}$ are each independently selected from the group consisting of hydrogen, $C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yl, $C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, Ar, Ar-$C_{1-6}$-

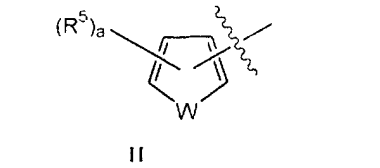
alk(en/yn)yl, Ar-$C_{3-8}$-cycloalk(en)yl, Ar-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, hydroxy-$C_{1-6}$-alk(en/yn)yl, hydroxy-$C_{3-8}$-cycloalk(en)yl, hydroxy-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, halo-$C_{1-6}$-alk(en/yn)yl, halo-$C_{3-8}$-cycloalk(en)yl, halo-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, cyano-$C_{1-6}$-alk(en/yn)yl, cyano-$C_{3-8}$-cycloalk(en)yl and cyano-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl[[,]]; or

$\mathbf{R^{12}}$ and $\mathbf{R^{12'}}$ taken together with the nitrogen atom to which they are attached form a 4-8 membered saturated or unsaturated ring [[which]]that optionally contains 1, 2 or 3 further heteroatoms;

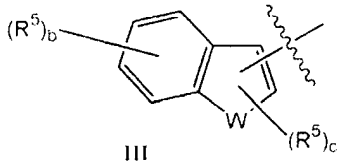
with the proviso that when $\mathbf{R^3}$ is $N\mathbf{R^{12}R^{12'}}$, then **q** is 0;

and

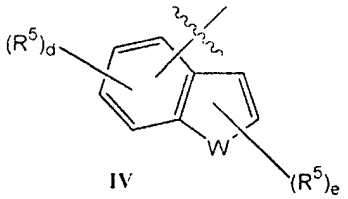
**Y** represents a group of formula **II**, **III**, **IV**, **V**, **VI**, **XXX** or **XXXI**:



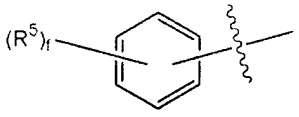

II



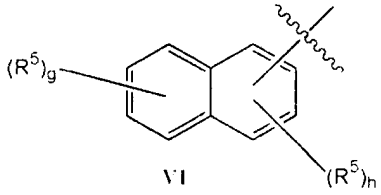

III



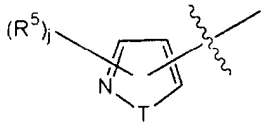

IV




V

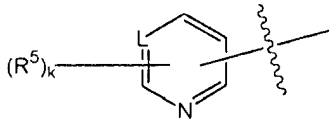



VI




XXX

or




XXXI

wherein:

**W** is O or S;

**T** is N, NH or O;

**L** is N, C or CH;

**a** is 0, 1, 2 or 3;

**b** is 0, 1, 2, 3 or 4;

**c** is 0 or 1;

**d** is 0, 1, 2 or 3;

**e** is 0, 1 or 2;

**f** is 0, 1, 2, 3, 4 or 5;

**g** is 0, 1, 2, 3 or 4;

**h** is 0, 1, 2 or 3;

**j** is 0, 1, 2 or 3; with the proviso that when **T** is a nitrogen atom, then **j** is 0, 1, 2 or 3; and when **T** is NH or an oxygen atom then **j** is 0, 1 or 2;

**k** is 0, 1, 2, 3 or 4; and

each $\mathbf{R^5}$ is independently selected from the group consisting of a $C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yl, $C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, Ar, Ar-$C_{1-6}$-alk(en/yn)yl, Ar-thio, Ar-oxy, acyl, $C_{1-6}$-alk(en/yn)yloxy, $C_{3-8}$-cycloalk(en)yloxy, $C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yloxy, halogen, halo-$C_{1-6}$-alk(en/yn)yl, halo-$C_{3-8}$-cycloalk(en)yl, halo-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, -CO-N$\mathbf{R^6R^{6'}}$, cyano, cyano-$C_{1-6}$-alk(en/yn)yl, cyano-$C_{3-8}$-cycloalk(en)yl, cyano-$C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, -N$\mathbf{R^7R^{7'}}$, -S-$\mathbf{R^8}$ and -

$SO_2R^8$[[,]]; or two adjacent $R^5$ groups taken together with the aromatic group to which they are attached form a 4-8 membered ring [[which]]that optionally contains one or two heteroatoms; wherein:

$R^6$ and $R^{6'}$ are each independently selected from the group consisting of hydrogen, $C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yl, $C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl and Ar;

$R^7$ and $R^{7'}$ are each independently selected from the group consisting of hydrogen, $C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yl, $C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, Ar and acyl;

and

$R^8$ is selected from the group consisting of hydrogen, $C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yl, $C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl, Ar and $-NR^9R^{9'}$; wherein:

$R^9$ and $R^{9'}$ are each independently selected from the group consisting of hydrogen, $C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yl and $C_{3-8}$-cycloalk(en)yl-$C_{1-6}$-alk(en/yn)yl; provided that when $R^8$ is $-NR^9R^{9'}$, then $R^5$ is not -S-$R^8$;

or salts thereof;

with the proviso that the compound of formula I is not:

N-[1-(phenylmethyl)-1H-indol-5-yl]-Methanesulfonamide;

N-[1-[(4-fluorophenyl)methyl]-1H-indol-5-yl]-Methanesulfonamide;

N-[2,3-dihydro-1-(phenylmethyl)-1H-indol-5-yl]-Methanesulfonamide;

N-[1-(phenylmethyl)-1H-indol-5-yl]-N'-4-quinolinyl-Urea;

N-[1-(phenylmethyl)-1H-indol-5-yl]-N'-4-quinolinyl-Urea; or

1-(1-benzyl-5-indolinyl)-3-phenyl-Urea;

or salts thereof.

2. (Original) A compound according to Claim 1, wherein at least one of $\mathbf{R^1}$ or $\mathbf{R^{1'}}$ is a hydrogen atom.

3. (Previously presented) A compound according to claim 2, wherein both $\mathbf{R^1}$ and $\mathbf{R^{1'}}$ are hydrogen atoms.

4. (Previously presented) A compound according to claim 1, wherein **s** is 0.

5. (Previously presented) A compound according to claim 1, wherein **s** is 1.

6. (Previously presented) A compound according to claim 1, wherein $\mathbf{R^2}$ is a hydrogen atom.

7. (Previously presented) A compound according claim 1, wherein $\mathbf{R^2}$ is $NO_2$ or a halogen atom.

8. (Previously presented) A compound according to claim 1, wherein **U** is $NR^{11}$.

9. (Currently Amended) A compound according to [[Claim]]claim 8, wherein $\mathbf{R^{11}}$ is a hydrogen atom.

10. (Previously presented) A compound according to claim 1, wherein **X** is CO.

11. (Previously presented) A compound according to claim 1, wherein **X** is $SO_2$.

12. (Previously presented) A compound according to claim 1, wherein **q** is 0.

13. (Previously presented) A compound according to claim 1, wherein **q** is 1.

14. (Currently amended) A compound according to [[Claim]]claim 13, wherein **Z** is an oxygen atom.

15. (Currently amended) A compound according to claim 1, wherein $\mathbf{R^3}$ is selected from the group consisting of $C_{1-6}$-alk(en/yn)yl, $C_{3-8}$-cycloalk(en)yl, Ar, Ar-$C_{1-6}$-alk(en/yn)yl, Ar-oxy-$C_{1-6}$-alk(en/yn)yl, Ar-$C_{1-6}$-alk(en/yn)yloxy-$C_{1-6}$-alk(en/yn)yl and -$N\mathbf{R^{12}R^{12'}}$; with the proviso that when $\mathbf{R^3}$ is $N\mathbf{R^{12}R^{12'}}$, then **q** is 0.

16. (Currently amended) A compound according to [[Claim]]claim 15, wherein $\mathbf{R^3}$ is $N\mathbf{R^{12}R^{12'}}$, **q** is 0 and $\mathbf{R^{12}}$ and $\mathbf{R^{12'}}$ are each independently selected from the group consisting of hydrogen, $C_{1-6}$-alk(en/yn)yl, Ar and Ar-$C_{1-6}$-alk(en/yn)yl[[,]]; or $\mathbf{R^{12}}$ and $\mathbf{R^{12'}}$ taken together

with the nitrogen atom to which they are attached form a 4-8 membered saturated or unsaturated ring [[which]]that optionally contains 1, 2 or 3 further heteroatoms.

17. (Previously presented) A compound according to claim 1, wherein **Y** is of formula **II**, **III**, **V**, **XXX**, **or XXXI**.

18. (Previously presented) A compound according to claim 17, wherein **Y** is of formula **II** or **III** and **W** is a sulphur atom.

19. (Previously presented) A compound according to claim 17, wherein **Y** is of formula **XXX** and **T** is a nitrogen atom or an oxygen atom.

20. (Previously presented) A compound according to claim 17, wherein **Y** is of formula **XXXI** and **L** is C or CH.

21. (Currently amended) A compound according to claim 1, wherein each $\mathbf{R^5}$ is independently selected from the group consisting of $C_{1-6}$-alk(en/yn)yl, Ar, Ar-thio, Ar-oxy, halogen and halo-$C_{1-6}$-alk(en/yn)yl; or two adjacent $\mathbf{R^5}$ taken together with the aromatic group to which they are attached form a 4-8 membered ring [[which]]that optionally contains one or two heteroatoms.

22. (Currently amended) A compound selected from the group consisting of:

N-[4-Chloro-1-(4-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

N-[4-Chloro-1-(5-chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

[1-(4-Fluorobenzyl)-2,3-dihydro-1H-indol-5-yl]-carbamic acid propyl ester[[;]];

N-[1-(4-Fluorobenzyl)-2,3-dihydro-1H-indol-5-yl]-C-phenyl-methanesulfonamide[[;]];

4-Fluoro-N-[1-(4-fluorobenzyl)-2,3-dihydro-1H-indol-5-yl]-benzamide[[;]];

N-[1-(4-Fluorobenzyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide[[;]];

N-[1-(4-Fluorobenzyl)-2,3-dihydro-1H-indol-5-yl]-2-thiophen-2-ylacetamide[[;]];

N-[1-(4-Fluorobenzyl)-2,3-dihydro-1H-indol-5-yl]-2-(4-fluorophenyl)-acetamide[[;]];

3-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-1,1-diisopropylurea[[;]];

Morpholine-4-carboxylic acid [1-(5-chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-amide[[;]];

Pyrrolidine-1-carboxylic acid [1-(5-chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-amide[[;]];

[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-carbamic acid 2-benzyloxyethyl ester[[;]];

3-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-1-methyl-1-propylurea[[;]];

[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-carbamic acid tert-butyl ester[[;]];

N-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-C-phenyl-methanesulfonamide[[;]];

Butane-1-sulfonic acid [1-(5-chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-amide[[;]];

N-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-4-fluorobenzamide[[;]];

N-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-2,2-dimethylpropionamide[[;]];

N-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-2-phenoxyacetamide[[;]];

N-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide[[;]];

N-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-butyramide[[;]];

Cyclopentanecarboxylic acid [1-(5-chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-amide[[;]];

N-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-2-thiophen-2-ylacetamide[[;]];

N-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-isonicotinamide[[;]];

N-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-4-dimethylaminobenzamide[[;]];

N-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-2-(4-fluorophenyl)-acetamide[[;]];

N-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-6-trifluoromethylnicotinamide[[;]];

1-tert-Butyl-3-[1-(5-chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-urea[[;]];

1-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-3-ethylurea[[;]];

1-Benzyl-3-[1-(5-chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-urea[[;]];

1-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-3-phenethylurea[[;]];

1-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-3-thiophen-2-ylurea[[;]];

1-[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-3-thiophen-3-ylurea[[;]];

[1-(5-Chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-carbamic acid propyl ester;

2,2-Dimethyl-N-[6-nitro-1-(4-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-propionamide[[;]];

N-[1-(5-Chlorothiophen-2-ylmethyl)-6-nitro-2,3-dihydro-1H-indol-5-yl]-2,2-dimethylpropionamide[[;]];

2-(4-Fluorophenyl)-N-[6-nitro-1-(4-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-acetamide[[;]];

N-[1-(5-Chlorothiophen-2-ylmethyl)-6-nitro-2,3-dihydro-1H-indol-5-yl]-2-(4-fluorophenyl)-acetamide;

N-[1-(5-Chlorothiophen-2-ylmethyl)-6-nitro-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

N-[6-Amino-1-(5-chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide[[;]];

N-[6-Amino-1-(4-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-2,2-dimethylpropionamide; [[;]];

N-[6-Amino-1-(5-chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-2,2-dimethylpropionamide[[;]];

N-[6-Amino-1-(4-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-2-(4-fluorophenyl)-acetamide[[;]];

N-[6-Amino-1-(4-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

N-[6-Amino-1-(4-fluorobenzyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

N-[6-Amino-1-(3-fluoro-4-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

N-[1-(5-Chlorothiophen-2-ylmethyl)-1H-indol-5-yl]-3,3-dimethylbutyramide;

N-[6-Bromo-1-(4-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

N-[6-Bromo-1-(5-chlorothiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

N-[1-(4-Chlorobenzyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

3,3-Dimethyl-N-[1-(4-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-butyramide;

N-[1-(4-Isopropylbenzyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

N-[1-(3-Fluoro-4-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

N-[1-(6-Chlorobenzo[1,3]dioxol-5-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

N-[1-(3,5-Dimethyl-1-phenyl-1H-pyrazol-4-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

N-[1-(2-Chloro-5-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

N-{1-[5-(4-Chlorophenoxy)-1,3-dimethyl-1H-pyrazol-4-ylmethyl]-2,3-dihydro-1H-indol-5-yl}-3,3-dimethylbutyramide;

3,3-Dimethyl-N-[1-(6-p-tolyloxy-pyridin-3-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-butyramide;

N-{1-[6-(4-Chlorophenylsulfanyl)-pyridin-3-ylmethyl]-2,3-dihydro-1H-indol-5-yl}-3,3-dimethylbutyramide;

N-{1-[6-(4-Cyanophenoxy)-pyridin-3-ylmethyl]-2,3-dihydro-1H-indol-5-yl}-3,3-dimethylbutyramide;

3,3-Dimethyl-N-[1-(6-trifluoromethylpyridin-3-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-butyramide;

3,3-Dimethyl-N-[1-(3-methyl-benzo[b]thiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-butyramide;

N-[1-(6-Fluoro-4H-benzo[1,3]dioxin-8-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-3,3-dimethylbutyramide;

3,3-Dimethyl-N-[1-(6-phenoxypyridin-3-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-butyramide;

3,3-Dimethyl-N-[1-(3-methyl-5-phenyl-isoxazol-4-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-butyramide;

N-(1-Benzo[b]thiophen-2-ylmethyl-2,3-dihydro-1H-indol-5-yl)-3,3-dimethylbutyramide;

N-{1-[1-(4-Fluorophenyl)-5-methyl-1H-pyrazol-4-ylmethyl]-2,3-dihydro-1H-indol-5-yl}-3,3-dimethylbutyramide;

3,3-Dimethyl-N-[1-(5-methylthiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-butyramide;

3,3-Dimethyl-N-[1-(4-pyrrol-1-yl-benzyl)-2,3-dihydro-1H-indol-5-yl]-butyramide;

N-[1-(4-Chlorobenzyl)-2,3-dihydro-1H-indol-5-yl]-2-(4-fluorophenyl)-acetamide;

2-(4-Fluorophenyl)-N-[1-(4-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-acetamide;

2-(4-Fluorophenyl)-N-[1-(4-isopropylbenzyl)-2,3-dihydro-1H-indol-5-yl]-acetamide;

2-(4-Fluorophenyl)-N-[1-(3-fluoro-4-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-acetamide;

N-[1-(6-Chlorobenzo[1,3]dioxol-5-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-2-(4-fluorophenyl)-acetamide;

N-[1-(3,5-Dimethyl-1-phenyl-1H-pyrazol-4-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-2-(4-fluorophenyl)-acetamide;

N-[1-(2-Chloro-5-trifluoromethylbenzyl)-2,3-dihydro-1H-indol-5-yl]-2-(4-fluorophenyl)-acetamide;

N-{1-[5-(4-Chlorophenoxy)-1,3-dimethyl-1H-pyrazol-4-ylmethyl]-2,3-dihydro-1H-indol-5-yl}-2-(4-fluorophenyl)-acetamide;

N-{1-[6-(4-Cyanophenoxy)-pyridin-3-ylmethyl]-2,3-dihydro-1H-indol-5-yl}-2-(4-fluorophenyl)-acetamide;

2-(4-Fluorophenyl)-N-[1-(3-methyl-benzo[b]thiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-acetamide;

N-[1-(6-Fluoro-4H-benzo[1,3]dioxin-8-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-2-(4-fluorophenyl)-acetamide;

2-(4-Fluorophenyl)-N-[1-(6-phenoxypyridin-3-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-acetamide;

N-(1-Benzo[b]thiophen-2-ylmethyl-2,3-dihydro-1H-indol-5-yl)-2-(4-fluorophenyl)-acetamide;

2-(4-Fluorophenyl)-N-{1-[1-(4-fluorophenyl)-5-methyl-1H-pyrazol-4-ylmethyl]-2,3-dihydro-1H-indol-5-yl}-acetamide;

2-(4-Fluorophenyl)-N-[1-(5-methylthiophen-2-ylmethyl)-2,3-dihydro-1H-indol-5-yl]-acetamide; and

2-(4-Fluorophenyl)-N-[1-(4-pyrrol-1-yl-benzyl)-2,3-dihydro-1H-indol-5-yl]-acetamide[[,]]; or

a pharmaceutically acceptable salt thereof.

23. (Previously presented) A pharmaceutical composition comprising a compound according to claim 1 and one or more pharmaceutically acceptable carriers or diluents.

24-37. (Withdrawn).